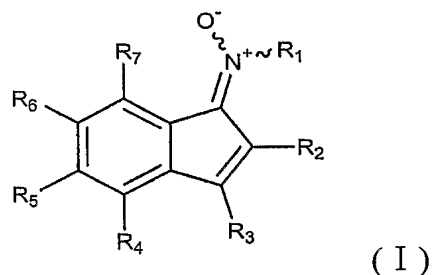


What is claimed is:

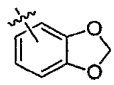

1. An indene derivative of formula (I) or a pharmaceutically acceptable salt thereof:



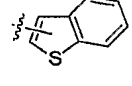



wherein,

$R_1$  is  $C_{1-6}$  alkyl,  $C_{1-6}$  alkenyl or  $C_{3-6}$  cycloalkyl, which is unsubstituted or substituted with one or more phenyl groups;

$R_2$  is H, CN,  $CO_2R^a$ ,  $CH_2CO_2R^a$ ,  $CONR^bR^c$ , , or phenyl;

$R_3$  is  $C_{1-6}$  alkyl,  $C_{3-6}$  cycloalkyl, or naphthyl, phenyl, , ,

, ,  or , which is unsubstituted or substituted with one or more substituents selected from the group consisting of halogen, CN,  $NH_2$ ,  $NO_2$ ,  $OR^a$ , phenyloxy,  $C_{1-6}$  alkyl and  $C_{3-6}$  cycloalkyl; and

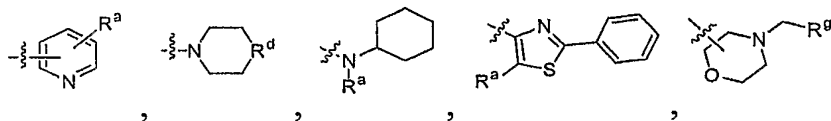
$R_4$ ,  $R_5$ ,  $R_6$  and  $R_7$  are each independently H, OH,  $OSO_2CH_3$ ,  $O(CH_2)_mR^e$ ,  $CH_2R^f$ ,  $OCOCH_2OR^g$ ,  $OCH_2CH_2OR^g$  or  $OCH_2CH=CHR^g$ , or  $R_5$  and  $R_6$  together form  $OCH_2O$ ;

in which  $R^a$  is H, or  $C_{1-6}$  alkyl or  $C_{3-6}$  cycloalkyl, which is unsubstituted or substituted with one or more halogens;

$R^b$  and  $R^c$  are each independently H,  $C_{1-6}$  alkyl or  $C_{3-6}$  cycloalkyl;

$R^d$  is O, S or  $NR^a$ ;

$R^e$  is H, halogen,  $C_{3-6}$  cycloalkyl, naphthyl,



or phenyl, which is

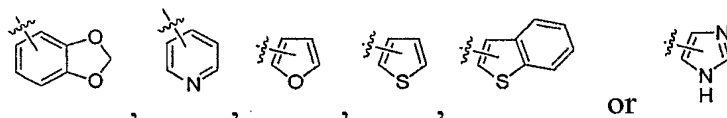
unsubstituted or substituted with one or more substituents selected from the group consisting of halogen, CN,  $NH_2$ ,  $NO_2$ ,  $OR^a$ ,  $CF_3$  and  $COOR^a$ ;

$R^f$  is  $OCH_2CH_2R^g$  or ;

$R^g$  is phenyl, which is unsubstituted or substituted with one or more substituents selected from the group consisting of halogen, CN,  $NH_2$ ,  $NO_2$  and  $OR^a$ ; and

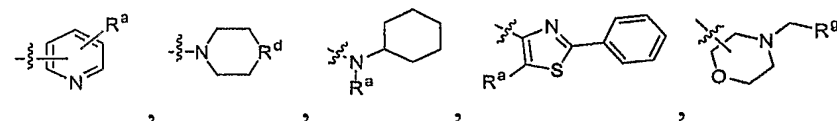
$m$  is an integer in the range of 1 to 5.

2. The compound of claim 1, wherein  $R_1$  is  $C_{1-6}$  alkyl, which is unsubstituted or substituted with a phenyl group;  $R_2$  is H, CN,  $CO_2R^a$ ,  $CH_2CO_2R^a$ ,  $CONR^bR^c$  or phenyl;  $R_3$  is  $C_{1-6}$  alkyl,  $C_{3-6}$  cycloalkyl, or phenyl,



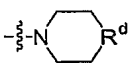
or , which is unsubstituted or

substituted with one or more substituents selected from the group consisting of halogen,  $C_{1-6}$  alkyl and  $C_{3-6}$  cycloalkyl;  $R_4$  and  $R_7$  are H;  $R_5$  and  $R_6$  are each independently OH,  $OSO_2CH_3$ ,  $O(CH_2)_mR^e$ ,  $CH_2R^f$ ,  $OCOCH_2OR^g$ ,  $OCH_2CH_2OR^g$  or  $OCH_2CH=CHR^g$ , or together form  $OCH_2O$ ;  $R^a$  is H, or  $C_{1-6}$  alkyl;  $R^d$  is O or  $NCH_3$ ;  $R^e$  is H, halogen,  $C_{3-6}$  cycloalkyl, naphthyl,

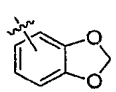



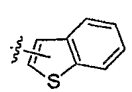



or phenyl, which is

unsubstituted or substituted with one or more substituents selected from the group consisting of halogen, OH, methoxy,  $CF_3$  and  $COOR^a$ ;  $R^f$  is  $OCH_2CH_2R^g$

or ; and R<sup>g</sup> is phenyl.

3. The compound of claim 2, wherein R<sub>1</sub> is CH<sub>3</sub>; R<sub>2</sub> is H, CN, CO<sub>2</sub>R<sup>a</sup> or

CONR<sup>b</sup>R<sup>c</sup>; R<sub>3</sub> is C<sub>1-6</sub> alkyl, or phenyl, , , , , 

5 or , which is unsubstituted or substituted with one or more halogens or C<sub>1-6</sub> alkyl groups; and R<sub>5</sub> and R<sub>6</sub> are each independently O(CH<sub>2</sub>)<sub>m</sub>R<sup>e</sup> or CH<sub>2</sub>R<sup>f</sup>, or together form OCH<sub>2</sub>O.

10 4. The compound of claim 1, which is selected from the group consisting of:

- 1) 6-methoxy-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 2) 1-(*trans*-isopropylimino-*N*-oxy)-6-methoxy-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 15 3) 1-(*trans*-benzylimino-*N*-oxy)-6-methoxy-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 4) 1-(*trans*-ethylimino-*N*-oxy)-6-methoxy-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 5) 6-methoxy-1-(*trans*-phenylpropylimino-*N*-oxy)-3-phenyl-1H-indene-2-
- 20 carboxylate ethyl ester
- 6) 6-methoxy-1-(*trans*-(2-methylbutenylimino)-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 7) 1-(*trans*-isobutylimino-*N*-oxy)-6-methoxy-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 25 8) 1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester

- 9) 1-(*trans*-methyylimino-*N*-oxy)-3-phenyl-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 10) 1-(*trans*-methyylimino-*N*-oxy)-6-phenetyloxy-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 5 11) 3-furan-3-yl-1-(*trans*-methyylimino-*N*-oxy)-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 12) 6-hydroxy-1-(*trans*-methyylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 13) 1-(*cis*-methyylimino-*N*-oxy)-3-phenyl-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 10 14) 3-(*trans*-methyylimino-*N*-oxy)-1-phenyl-3H-indene-5-ol
- 15) 1-(*trans*-methyylimino-*N*-oxy)-3-phenyl-6-(5-phenylpentyloxy)-1H-indene-2-carboxylate ethyl ester
- 16) 1-(*cis*-methyylimino-*N*-oxy)-3-phenyl-6-(5-phenylpentyloxy)-1H-indene-2-carboxylate ethyl ester
- 15 17) 6-[2-(4-chlorophenoxy)acetoxyl]-1-(*trans*-methyylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 18) 6-[2-(4-chlorophenoxy)ethoxyl]-1-(*trans*-methyylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 20 19) 1-(*trans*-methyylimino-*N*-oxy)-6-(naphthalene-2-ylmethoxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 20) methyl-[3-phenyl-6-(3-phenylpropoxy)indene-1-ylidene]amine-*N*-oxide
- 21) 1-(*trans*-methyylimino-*N*-oxy)-6-[2-(5-methyl-2-phenylthiazol-4-yl)ethoxyl]-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 25 22) 1-(*trans*-methyylimino-*N*-oxy)-3-phenyl-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 23) 6-[2-(4-hydroxyphenyl)ethoxyl]-1-(*trans*-methyylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 24) 6-(2-adaman-1-ylethoxy)-1-(*trans*-methyylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 30 25) 6-(2-cyclohexylethoxy)-1-(*trans*-methyylimino-*N*-oxy)-3-phenyl-1H-indene-

2-carboxylate ethyl ester

26) 1-(*trans*-methyylimino-*N*-oxy)-3-phenyl-6-(3-phenylpropenoxy)-1H-indene-2-carboxylate ethyl ester

27) 6-[2-(2-fluorophenyl)ethoxy]-1-(*trans*-methyylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester

28) 6-[2-(3-fluorophenyl)ethoxy]-1-(*trans*-methyylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester

29) 6-[2-(4-fluorophenyl)ethoxy]-1-(*trans*-methyylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester

30) 1-(*trans*-methyylimino-*N*-oxy)-3-phenyl-6-[2-(3-trifluoromethylphenyl)ethoxy]-1H-indene-2-carboxylate ethyl ester

31) 6-(4-methoxycarbonylbenzyloxy)-1-(*trans*-methyylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester

32) 1-(*trans*-methyylimino-*N*-oxy)-3-phenyl-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl amide

33) 1-(*trans*-methyylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester

34) 6-[2-(cyclohexylmethylamino)ethoxy]-1-(*trans*-methyylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester

35) 3-(2-fluorophenyl)-6-methoxy-1-(*trans*-methyylimino-*N*-oxy)-1H-indene-2-carboxylate ethyl ester

36) 1-(*trans*-methyylimino-*N*-oxy)-6-[2-(4-methylpiperazine-1-yl)ethoxy]-3-phenyl-1H-indene-2-carboxylate ethyl ester

37) (2,3-diphenyl indene-1-yl lidene)methylamine-*N*-oxide

38) 1-(*trans*-methyylimino-*N*-oxy)-3-phenyl-6-(3-phenylpropoxy)-1H-indene-2-carboxylate isopropyl amide

39) 1-(*trans*-methyylimino-*N*-oxy)-3-phenyl-6-(3-phenylpropoxy)-1H-indene-2-carboxylate cyclohexyl amide

40) [1-(*trans*-methyylimino-*N*-oxy)-3-phenyl-6-(3-phenylpropoxy)-1H-indene-2-yl]morpholine-4-yl-methanone

41) 1-(*trans*-methyylimino-*N*-oxy)-6-(2-morpholine-4-yl-ethoxy)-3-phenyl-1H-

indene-2-carboxylate cyclohexyl amide

42) 1-(*trans*-methylimino-*N*-oxy)-3-phenyl-5-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester

43) 1-(*trans*-methylimino-*N*-oxy)-6-phenethyloxymethyl-3-phenyl-1H-indene-2-carboxylate ethyl ester

44) (6-methoxy-3-phenylindene-1-ylidene)methylamine-*N*-oxide

45) 1-(*cis*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester

46) 6-(2-bromoethoxy)-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester

47) 1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-3-phenyl-1H-indene-2-carboxylate *tert*-buthyl ester

48) 1-(*trans*-methylimino-*N*-oxy)-5,6-methylenedioxy-1-oxo-3-phenyl-1H-indene-2-carboxylate ethyl ester

49) 4-[2-isopropylcarbamoyl-3-(*trans*-methylimino-*N*-oxy)-1-phenyl-3H-indene-5-yl-oxylmethyl]benzoate methyl ester

50) 1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-3-phenyl-1H-indene-2-carboxylate isopropyl amide

51) 1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-3-phenyl-1H-indene-2-carboxylate cyclopropyl amide

52) 3-(3-fluorophenyl)-1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-1H-indene-2-carboxylate isopropyl amide

53) (6-methoxy-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-yl)acetate ethyl ester

54) (6-methoxy-1-(*cis*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-yl)acetate ethyl ester

55) 5-[2-(5-ethylpyridine-2-yl)ethoxy]-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate isopropyl amide

56) 1-(*trans*-methylimino-*N*-oxy)-6-(3-phenylpropoxy)-3-*p*-tolyl-1H-indene-2-carboxylate ethyl ester

57) 1-(*trans*-methylimino-*N*-oxy)-6-(3-phenylpropoxy)-3-thiophene-2-yl-1H-

- indene-2-carboxylate ethyl ester
- 58) 3-(4-chlorophenyl)-1-(*trans*-methyylimino-*N*-oxy)-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 59) 3-(5-chlorothiophene-2-yl)-1-(*trans*-methyylimino-*N*-oxy)-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 60) 1-(*trans*-methyylimino-*N*-oxy)-6-(3-phenylpropoxy)-3-*m*-tolyl-1H-indene-2-carboxylate ethyl ester
- 61) 1-(*trans*-methyylimino-*N*-oxy)-3-(4-phenoxyphenyl)-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 62) 3-benzo-[1,3]-dioxol-5-yl-1-(*trans*-methyylimino-*N*-oxy)-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 63) methyl-[6-(3-phenylpropoxy)-3-pyridine-2-yl-indene-1-ylidene]-amine-*N*-oxide
- 64) 3-furan-2-yl-1-(*trans*-methyylimino-*N*-oxy)-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 65) 3-ethyl-1-(*trans*-methyylimino-*N*-oxy)-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 66) 3-methyl-1-(*trans*-methyylimino-*N*-oxy)-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 67) 1-(*trans*-methyylimino-*N*-oxy)-6-(3-phenylpropoxy)-3-thiophene-3-yl-1H-indene-2-carboxylate ethyl ester
- 68) 3-cyclopropyl-1-(*trans*-methyylimino-*N*-oxy)-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 69) 1-(*trans*-methyylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-3-thiophene-3-yl-1H-indene-2-carboxylate ethyl ester
- 70) 3-benzo-[b]-thiophene-3-yl-1-(*trans*-methyylimino-*N*-oxy)-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 71) 3-(1H-imidazole-4-yl)-1-(*trans*-methyylimino-*N*-oxy)-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 72) 3-(1-ethyl propyl)-1-(*trans*-methyylimino-*N*-oxy)-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester

- 73) 1-(*trans*-methyylimino-*N*-oxy)-3-phenyl-6-(3-phenylpropoxy)-1H-indene-2-carboxylate amide
- 74) 6-(4-benzylmorpholine-2-ylmethoxy)-1-(*trans*-methyylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate isopropyl amide
- 5 75) 1-(*trans*-methyylimino-*N*-oxy)-3-phenyl-6-(3-phenylpropoxy)-1H-indene-2-carbonitrile
- 76) 1-(*trans*-methyylimino-*N*-oxy)-5,6-methylenedioxy-1-oxo-3-phenyl-1H-phenyl-2-carboxylate isopropyl amide
- 77) 1-(*trans*-methyylimino-*N*-oxy)-6-morpholine-4-ylmethyl-3-phenyl-1H-  
10 indene-2-carboxylate ethyl ester
- 78) 1-(*trans*-methyylimino-*N*-oxy)-3-phenyl-6-(2-pyridine-2-ylethoxy)-1H-indene-2-carboxylate ethyl ester
- 79) 6-[2-(5-ethylpyridine-2-yl)ethoxy]-1-(*trans*-methyylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 15 80) 1-(*trans*-methyylimino-*N*-oxy)-3-phenyl-6-(2-pyridine-2-ylethoxy)-1H-indene-2-carboxylate isopropyl amide
- 81) 6-[2-(5-ethylpyridine-2-yl)ethoxy]-1-(*trans*-methyylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate isopropyl amide
- 82) methyl-[6-(2-morpholine-4-ylethoxy)-3-phenylindene-1-ylidene]amine-*N*-  
20 oxide
- 83) 5,6-bis-methanesulfonyloxy-1-(*trans*-methyylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 84) 1-(*trans*-methyylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-3-phenyl-1H-indene-2-carboxylate isobutyl ester
- 25 85) 1-(*trans*-methyylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-3-phenyl-1H-indene-2-carboxylate methyl ester
- 86) 1-(*cis*-methyylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-3-phenyl-1H-indene-2-carboxylate methyl ester
- 87) 1-(*trans*-methyylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-3-phenyl-1H-  
30 indene-2-carboxylate propyl ester
- 88) 3-(4-fluorophenyl)-1-(*trans*-methyylimino-*N*-oxy)-6-(2-morpholine-4-



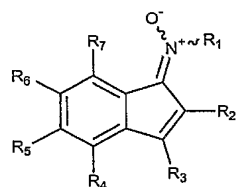
- ylethoxy)-1H-indene-2-carboxylate ethyl ester
- 89) 1-(*trans*-methylimino-*N*-oxy)-3-phenyl-6-(pyridine-2-ylmethoxy)-1H-indene-2-carboxylate ethyl ester
- 90) 1-(*trans*-methylimino-*N*-oxy)-3-phenyl-6-(pyridine-2-yloxy)-1H-indene-2-carboxylate ethyl ester
- 5 91) 6-(3-methoxybenzyloxy)-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 92) 1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-3-thiophene-3-yl-1H-indene-2-carboxylate isopropyl amide
- 10 93) 3-(1-ethylpropyl)-1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-1H-indene-2-carboxylate ethyl ester
- 94) 3-benzo-[b]-thiophene-3-yl-1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-1H-indene-2-carboxylate isopropyl amide
- 95) 3-(4-fluorophenyl)-1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-1H-indene-2-carboxylate isopropyl amide
- 15 96) 3-(1-ethylpropyl)-1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-1H-indene-2-carboxylate isopropyl amide
- 97) 1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-3-(2,4,6-trimethylphenyl)-1H-indene-2-carboxylate ethyl ester
- 20 98) 3-(2,6-dimethylphenyl)-1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-1H-indene-2-carboxylate ethyl ester
- 99) 1-(*trans*-methylimino-*N*-oxy)-3-phenyl-5-(2-pyridine-2-ylethoxy)-1H-indene-2-carboxylate isopropyl amide
- 100) 1-(*trans*-methylimino-*N*-oxy)-5-(2-morpholine-4-ylethoxy)-3-phenyl-1H-indene-2-carboxylate isopropyl amide
- 25 101) 1-(*cis*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-3-phenyl-1H-indene-2-carboxylate isopropyl ester
- 102) 3-(3-fluorophenyl)-1-(*trans*-methylimino-*N*-oxy)-6-(2-pyridine-2-ylethoxy)-1H-indene-2-carboxylate isopropyl amide
- 30 103) 6-[2-(5-ethylpyridine-2-yl)ethoxy]-3-(3-fluorophenyl)-1-(*trans*-methylimino-*N*-oxy)-1H-indene-2-carboxylate isopropyl amide

104) 3-(4-cyanophenyl)-6-(2-morpholine-4-ylethoxy)-1-(*trans*-methylimino-*N*-oxy)-1H-indene-2-carboxylate ethyl ester

105) 1-(*trans*-methylimino-*N*-oxy)-3-phenyl-6-(2-pyridine-2-ylethoxy)-1H-indene-2-carboxylate isopropyl ester.

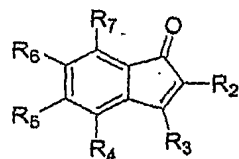
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5. A process for preparing the indene derivative of claim 1, which comprises step of subjecting indenone compound of formula (II) to a condensation reaction with  $R_1\text{NHOH}$  or  $\text{NH}_2\text{OH}$  to obtain a compound of formula (III), and reacting the compound of formula (III) with  $R_1\text{X}$ :

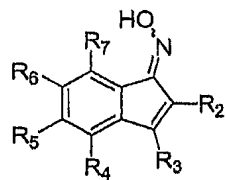


( I )

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( II )



( III )

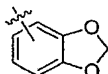

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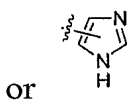
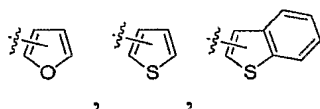
wherein,

X is halogen;

$R_1$  is  $\text{C}_{1-6}$  alkyl,  $\text{C}_{1-6}$  alkenyl or  $\text{C}_{3-6}$  cycloalkyl, which is unsubstituted or substituted with one or more phenyl groups;

$R_2$  is H, CN,  $\text{CO}_2R^a$ ,  $\text{CH}_2\text{CO}_2R^a$ ,  $\text{CONR}^bR^c$ , , or phenyl;

$R_3$  is  $\text{C}_{1-6}$  alkyl,  $\text{C}_{3-6}$  cycloalkyl, or naphthyl, phenyl, , ,



or which is unsubstituted or substituted with one or more substituents selected from the group consisting of halogen, CN,  $\text{NH}_2$ ,  $\text{NO}_2$ ,  $\text{OR}^a$ , phenyloxy,  $\text{C}_{1-6}$  alkyl and  $\text{C}_{3-6}$  cycloalkyl; and

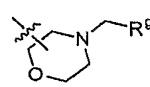
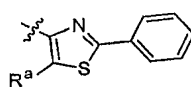
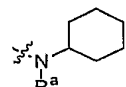
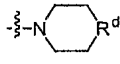
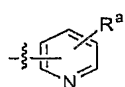
$R_4$ ,  $R_5$ ,  $R_6$  and  $R_7$  are each independently H, OH,  $\text{OSO}_2\text{CH}_3$ ,  $\text{O}(\text{CH}_2)_mR^e$ ,  $\text{CH}_2R^f$ ,  $\text{OCOCH}_2\text{OR}^g$ ,  $\text{OCH}_2\text{CH}_2\text{OR}^g$  or  $\text{OCH}_2\text{CH}=\text{CHR}^g$ , or  $R_5$  and  $R_6$  together form  $\text{OCH}_2\text{O}$ ;

in which  $R^a$  is H, or  $\text{C}_{1-6}$  alkyl or  $\text{C}_{3-6}$  cycloalkyl, which is unsubstituted or substituted with one or more halogens;

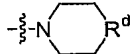
$R^b$  and  $R^c$  are each independently H,  $\text{C}_{1-6}$  alkyl or  $\text{C}_{3-6}$  cycloalkyl;

$R^d$  is O, S or  $\text{NR}^a$ ;

$R^e$  is H, halogen,  $\text{C}_{3-6}$  cycloalkyl, naphthyl,



or phenyl, which is unsubstituted or substituted with one or more substituents selected from the group consisting of halogen, CN,  $\text{NH}_2$ ,  $\text{NO}_2$ ,  $\text{OR}^a$ ,  $\text{CF}_3$  and  $\text{COOR}^a$ ;

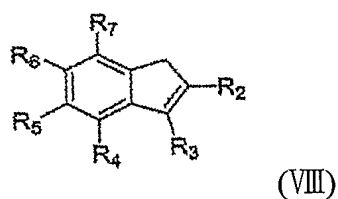
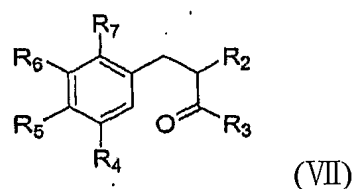
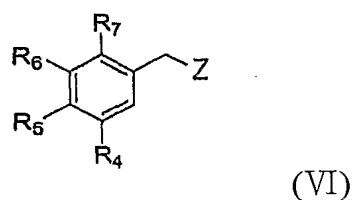
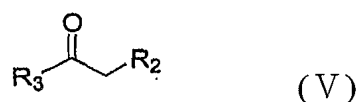
$R^f$  is  $\text{OCH}_2\text{CH}_2R^g$  or ;

$R^g$  is phenyl, which is unsubstituted or substituted with one or more substituents selected from the group consisting of halogen, CN,  $\text{NH}_2$ ,  $\text{NO}_2$  and  $\text{OR}^a$ ; and

$m$  is an integer in the range of 1 to 5.

6. The process of claim 5, wherein the indenone compound of formula (II) is prepared by a process comprising the steps of:

- 1) reacting compounds of formula (V) and (VI) to obtain a compound of formula (VII);
- 2) subjecting the compound of formula (VII) to cyclization to obtain a compound of formula (VIII); and
- 3) subjecting the compound of formula (VIII) to oxidation.

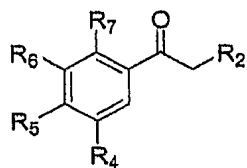


wherein,

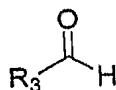
R<sub>2</sub> to R<sub>7</sub> have the same meanings as defined in claim 5, and Z is halogen or activated leaving group.

7. The process of claim 5, wherein the indenone compound of formula (II) is prepared by a process comprising the steps of:

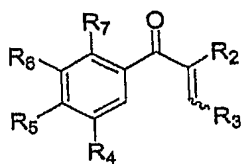
- 1) reacting compounds of formula (IX) and (X) to obtain a compound of formula (XI);
- 2) subjecting the compound of formula (XI) to cyclization to obtain a compound of formula (XII); and
- 3) subjecting the compound of formula (XII) to oxidation.



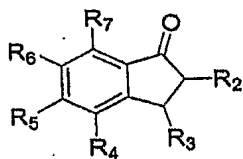
(IX)



(X)



(XI)



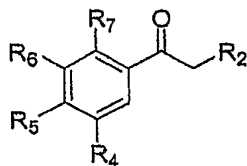
(XII)

wherein,

$R_2$  to  $R_7$  have the same meanings as defined in claim 5.

8. The process of claim 5, wherein the indenone compound of formula (II) is prepared by a process comprising the steps of:

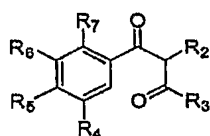
- 1) reacting compounds of formula (IX) and (XIII) to obtain a compound of formula (XIV); and
- 2) subjecting the compound of formula (XIV) to cyclization.



(IX)



(XIII)



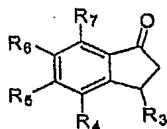
(XIV)

wherein,

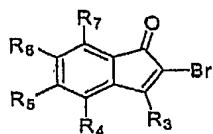
5  $R_2$  to  $R_7$  have the same meanings as defined in claim 5.

9. The process of claim 5, wherein the indenone compound of formula (II) is prepared by a process comprising the steps of:

- 10 1) subjecting a compound of formula (XV) to bromination obtain a compound of formula (XVI); and
- 2) subjecting the compound of formula (XVI) to a carbon-carbon coupling reaction in the presence of a metal catalyst, or to a substitution reaction using a nucleophile.



(XV)



(XVI)

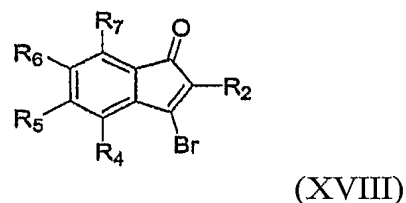
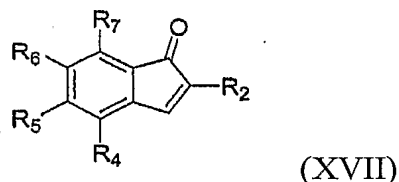
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wherein,

R<sub>3</sub> to R<sub>7</sub> have the same meanings as defined in claim 5.

10. The process of claim 5, wherein the indenone compound of formula (II) is prepared by a process comprising the steps of:

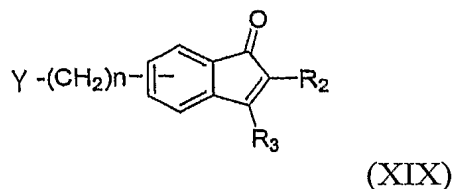
- 5 1) subjecting a compound of formula (XVII) to bromination obtain a compound of formula (XVIII); and
- 2) subjecting the compound of formula (XVIII) to a carbon-carbon coupling reaction in the presence of a metal catalyst, or to a substitution reaction using a nucleophile.



wherein,

R<sub>2</sub> and R<sub>4</sub> to R<sub>7</sub> have the same meanings as defined in claim 5.

15 11. The process of claim 5, wherein the indenone compound of formula (II) is prepared by subjecting a compound of formula (XIX) to an acylation reaction, a halogenation reaction followed by a substitution reaction by a nucleophile, or a carbon-carbon coupling reaction in the presence of a metal catalyst.



wherein,

$R_2$  and  $R_3$  have the same meanings as defined in claim 5, Y is hydroxy, thiol, amino  $C_{1-6}$  alkyl or halogen, and n is an integer in the range of 0 to 5.

5 12. A pharmaceutical composition for modulating the activities of peroxisome proliferator activated receptors (PPARs) comprising a therapeutically effective amount of the compound or salt defined in claim 1 as an active ingredient together with a pharmaceutically acceptable carrier.

10 13. The composition of claim 12, which is used for the treatment and prevention of diabetes, obesity, arteriosclerosis, hyperlipidemia, hyperinsulinism, hypertension, osteoporosis, liver cirrhosis, asthma and cancer.